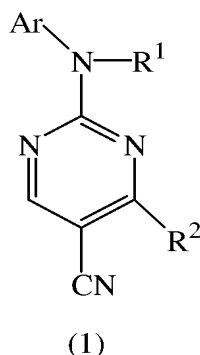


This listing of claims will replace all prior versions, and listings, of claims in the application.

**Listing of Claims:**

1. (currently amended) A compound of formula (1):



wherein

Ar is an optionally substituted aromatic or heteroaromatic group;

R<sup>1</sup> is a hydrogen atom or a straight or branched chain alkyl group;

R<sup>2</sup> is a -X<sup>1</sup>-R<sup>3</sup> group;

X<sup>1</sup> is a linker group selected from -C(O)-, -C(S)-, -S(O)-, -S(O)<sub>2</sub>-, -N(C<sub>1-6</sub> alkyl)-, -C(R<sup>7</sup>)<sub>2</sub>-, -CON(R<sup>7</sup>)-, -OC(O)N(R<sup>7</sup>)-, -CSN(R<sup>7</sup>)-, -N(R<sup>7</sup>)CO-, -N(R<sup>7</sup>)C(O)O-, -N(R<sup>7</sup>)CS-, -SON(R<sup>7</sup>)-, -SO<sub>2</sub>N(R<sup>7</sup>)-, -N(R<sup>7</sup>)SO<sub>2</sub>-, -N(R<sup>7</sup>)CON(R<sup>7</sup>)-, -N(R<sup>7</sup>)CSN(R<sup>7</sup>)-, -N(R<sup>7</sup>)SON(R<sup>7</sup>)-, and -N(R<sup>7</sup>)SO<sub>2</sub>N(R<sup>7</sup>)-;

R<sup>7</sup> is a hydrogen atom or C<sub>1-6</sub> alkyl group;

R<sup>3</sup> is an optionally substituted aliphatic, cycloaliphatic, heteroaliphatic, heterocycloaliphatic, aromatic or heteroaromatic group;

~~and the salts, solvates, hydrates and N-oxides~~ or a salt or N-oxide thereof.

2. (original) A compound according to Claim 1 wherein R<sup>1</sup> is a hydrogen atom.

- 3-5. (canceled)

6. (currently amended) A compound according to Claim 1 wherein Ar is a phenyl, pyridyl, indolyl, indazolyl, benzimidazolyl, benzothiazolyl, quinolyl, isoquinolyl or benzoxazolyl group each substituted by one, two or three -R<sup>4</sup> or -Alk(R<sup>4</sup>)<sub>m</sub> substituents;

~~in which~~ R<sup>4</sup> is a halogen atom, or an amino (-NH<sub>2</sub>), substituted amino, nitro, cyano, hydroxyl (-OH), substituted hydroxyl, formyl, carboxyl (-CO<sub>2</sub>H), esterified carboxyl, thiol (-SH), substituted thiol, -COR<sup>5</sup> [~~where R<sup>5</sup> is a -Alk(R<sup>4</sup>)<sub>m</sub>, aryl or heteroaryl group~~], -CSR<sup>5</sup>, -SO<sub>3</sub>H, -SO<sub>2</sub>R<sup>5</sup>, -SO<sub>2</sub>NH<sub>2</sub>, -SO<sub>2</sub>NHR<sup>5</sup>, ~~-SO<sub>2</sub>N[R<sup>5</sup>]<sub>2</sub>-SO<sub>2</sub>N(R<sup>5</sup>)<sub>2</sub>~~, -CONH<sub>2</sub>, -CSNH<sub>2</sub>, -CONHR<sup>5</sup>, -CSNHR<sup>5</sup>, ~~-CON[R<sup>5</sup>]<sub>2</sub>-CON(R<sup>5</sup>)<sub>2</sub>~~, ~~-CSN[R<sup>5</sup>]<sub>2</sub>-CSN(R<sup>5</sup>)<sub>2</sub>~~, -NHSO<sub>2</sub>H, -NHSO<sub>2</sub>R<sup>5</sup>, ~~-N[SO<sub>2</sub>R<sup>5</sup>]<sub>2</sub>-N(SO<sub>2</sub>R<sup>5</sup>)<sub>2</sub>~~, -NHSO<sub>2</sub>NH<sub>2</sub>, -NHSO<sub>2</sub>NHR<sup>5</sup>, ~~-NHSO<sub>2</sub>N[R<sup>5</sup>]<sub>2</sub>-NHSO<sub>2</sub>N(R<sup>5</sup>)<sub>2</sub>~~, -NHCOR<sup>5</sup>, -NHCONH<sub>2</sub>, -NHCONHR<sup>5</sup>, ~~-NHCON[R<sup>5</sup>]<sub>2</sub>-NHCON(R<sup>5</sup>)<sub>2</sub>~~, -NHCSR<sup>5</sup>, -NHC(O)OR<sup>5</sup>, or optionally substituted cycloaliphatic, hetero-cycloaliphatic, aryl or heteroaryl group;

R<sup>5</sup> is a -Alk(R<sup>4</sup>)<sub>m</sub>, aryl or heteroaryl group;

Alk is a straight or branched C<sub>1-6</sub> alkylene, C<sub>2-6</sub> alkenylene or C<sub>2-6</sub> alkynylene chain, optionally interrupted by one, two or three -O- or -S- atoms or groups selected from -S(O)-, -S(O)<sub>2</sub>- ~~or~~ and -N(R<sup>6</sup>)- [~~where R<sup>6</sup> is a hydrogen atom or a straight or branched chain C<sub>1-6</sub> alkyl group~~];

R<sup>6</sup> is a hydrogen atom or a straight or branched chain C<sub>1-6</sub> alkyl group; and

m is zero or an integer 1, 2 or 3.

7. (original) A compound according to Claim 6 wherein Ar is a phenyl group substituted by one, two or three -R<sup>4</sup> or -Alk(R<sup>4</sup>)<sub>m</sub> substituents.

8. (currently amended) A compound according to ~~any one of Claim 5 to Claim 7~~ claim 6 or claim 7 wherein at least one of -R<sup>4</sup>; and -Alk(R<sup>4</sup>)<sub>m</sub>; ~~R<sup>4b</sup> or Alk(R<sup>4b</sup>)<sub>m</sub>~~ is a -X<sup>1a</sup>(Alk<sup>a</sup>)<sub>p</sub>NR<sup>7a</sup>R<sup>7b</sup>) (~~where X<sup>1a</sup> is a direct bond or a linker atom or group, Alk<sup>a</sup> is as defined for Alk, p is zero or an integer 1 and R<sup>7a</sup> and R<sup>7b</sup> which may be the same or different is each a hydrogen atom or a straight or branched C<sub>1-6</sub> alkyl group~~), -X<sup>1a</sup>(Alk<sup>a</sup>)<sub>p</sub>NHet<sup>1</sup> (~~where NHet<sup>1</sup> is an optionally substituted C<sub>3-7</sub> cyclicamino group optionally containing one or more O or S~~

~~atoms or -N(R<sup>6</sup>) [where R<sup>6</sup> is a hydrogen atom or a straight or branched chain C<sub>1-6</sub>alkyl group] or -X<sup>1a</sup>(Alk<sup>a</sup>)<sub>p</sub>Ar<sup>2</sup> group (where Ar<sup>2</sup> is a nitrogen-containing heteroaromatic group);~~

X<sup>1a</sup> is a direct bond or a linker group selected from -C(O)-, -C(S)-, -S(O)-, -S(O)<sub>2</sub>-, -N(R<sup>7</sup>)-, -C(R<sup>7</sup>)<sub>2</sub>-, -CON(R<sup>7</sup>)-, -OC(O)N(R<sup>7</sup>)-, -CSN(R<sup>7</sup>)-, -N(R<sup>7</sup>)CO-, -N(R<sup>7</sup>)C(O)O-, -N(R<sup>7</sup>)CS-, -SON(R<sup>7</sup>)-, -SO<sub>2</sub>N(R<sup>7</sup>)-, -N(R<sup>7</sup>)SO<sub>2</sub>-, -N(R<sup>7</sup>)CON(R<sup>7</sup>)-, -N(R<sup>7</sup>)CSN(R<sup>7</sup>)-, -N(R<sup>7</sup>)SON(R<sup>7</sup>)-, and -N(R<sup>7</sup>)SO<sub>2</sub>N(R<sup>7</sup>)-;

Alk<sup>a</sup> is a straight or branched C<sub>1-6</sub> alkylene, C<sub>2-6</sub> alkenylene or C<sub>2-6</sub> alkynylene chain, optionally interrupted by one, two or three -O- or -S- atoms or groups selected from -S(O)-, -S(O)<sub>2</sub>- and -N(R<sup>6</sup>)-;

p is zero or an integer 1;

R<sup>7a</sup> and R<sup>7b</sup> are each independently a hydrogen atom or a straight or branched C<sub>1-6</sub>alkyl group;

-NHet<sup>1</sup> is an optionally substituted C<sub>3-7</sub>cyclicamino group optionally containing one or more -O- or -S- atoms or -N(R<sup>6</sup>); and

Ar<sup>2</sup> is a nitrogen containing heteroaromatic group.

9-10. (canceled)

11. (original) A pharmaceutical composition comprising a compound according to Claim 1 together with one or more pharmaceutically acceptable carriers, excipients or diluents.